

INFORMATION DISCLOSURE STATEMENT	Atty. Docket No.: 6315.N	Serial No.: 09/772,598
	Applicant(s): T. Benson et al.	Confirmation No.: 2967
	Filing Date: January 30, 2001	Group: 2183

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U.S. PATENT DOCUMENTS

Examiner Initial	Document Number	Date	Name	Class	Subclass	Technology Approval
	none					

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FOREIGN PATENT DOCUMENTS

Examiner Initial	Document Number	Date	Country	Class	Subclass	Translation	
						Yes	No
CSM	✓ WO 93/02209 A1	02/04/93	PCT				
	✓ WO 97/15588 A1	05/01/97	PCT				
	✓ EP 0786 519 A2	07/30/97	Europe				
	✓ WO 98/58961 A1	12/30/98	PCT				
	✓ WO 99/36422 A1	07/22/99	PCT				
	✓ WO 99/47639 A2	09/23/99	PCT				
	✓ WO 99/47662 A1	09/23/99	PCT				
	✓ WO 00/12678 A2	03/09/00	PCT				
	✓ WO 00/12678 A3	03/09/00	PCT				
	✓ WO 01/16292 A2	03/08/01	PCT				
✓							

OTHER DOCUMENTS (Including Authors, Title, Date, Pertinent Papers, etc.)

Examiner Initial	Document Description
CSM	✓ Bartlett et al., "CAVEAT: A program to facilitate the structure-derived design of biologically active molecules," <i>Molecular Recognition: Chemical and Biological Problems</i> , Royal Society of Chemistry, Special Pub No. 78:182-196 (1989).
↓	✓ Benson et al. "An enzyme-substrate complex involved in bacterial cell wall biosynthesis," <i>Nat Struct Biol.</i> 1995 Aug;2(8):644-53.
✓	✓ Blundell et al., <i>Protein Crystallography</i> , Academic Press, New York, NY; title page, publication page, and table of contents only, 8 pages (1976).

EXAMINER C. M. 17	Date Considered September 16, 2002
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Technology Center 2100

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✓	✓ Böhm, "The computer program LUDI: a new method for the de novo design of enzyme inhibitors," <i>J. Comput. Aided Mol. Des.</i> 1992 Feb;6(1):61-78.
✓	✓ Brünger, <i>X-plor Manual (Version 3.1) A System for X-ray Crystallography and NMR</i> (title page, publisher's page, and table of contents only), Yale University Press, New Haven, CT, 1992; 13 pgs.
✓	✓ Collaborative Computational Project, No. 4, "The CCP4 suite: programs for protein crystallography" <i>Acta Cryst.</i> 1994;D50:760-3.
✓	✓ Cowtan et al., "Improvement of Macromolecular Electron-Density Maps by the Simultaneous Application of Real and Reciprocal Space Constraints," <i>Acta Crystallogr D Biol Crystallogr.</i> 1993 Jan 1;49(Pt 1):148-157.
✓	✓ Cowtan et al., "Miscellaneous algorithms for density modification," <i>Acta Crystallogr D Biol Crystallogr.</i> 1998 Jul 1;54(Pt 4):487-93.
✓	✓ Eisen et al., "HOOK: a program for finding novel molecular architectures that satisfy the chemical and steric requirements of a macromolecule binding site," <i>Proteins: Struct. Funct. Genet.</i> 1994 Jul;19(3):199-221.
✓	✓ Evans, "SETOR: hardware-lighted three-dimensional solid model representations of macromolecules," <i>J Mol Graph.</i> 1993 Jun;11(2):134-8, 127-8.
✓	✓ Finzel, "LORE: exploiting database of known structures," <i>Meth. Enzymol.</i> 1997; 277(B):230-42.
✓	✓ Gillet et al., "SPROUT: a program for structure generation," <i>J. Comput. Aided Mol. Des.</i> 1993 Apr;7(2):127-53.
✓	✓ Goodford, "A computational procedure for determining energetically favorable binding sites on biologically important macromolecules," <i>J Med Chem.</i> 1985 Jul; 28(7):849-57.
✓	✓ Goodsell et al., "Automated docking of substrates to proteins by simulated annealing," <i>Proteins: Struct. Funct. Genet.</i> 1990;8(3):195-202.
✓	✓ Hendrickson, "Determination of macromolecular structures from anomalous diffraction of synchrotron radiation," <i>Science.</i> 1991 Oct 4;254(5028):51-8.
✓	✓ Hughes et al., "Structural gene for NAD synthetase in <i>Salmonella typhimurium</i> ," <i>J Bacteriol.</i> 1988 May;170(5):2113-20.

EXAMINER C.M. / 8	Date Considered September 16, 2002
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✓	Jiang et al., "Protein hydration observed by X-ray diffraction. Solvation properties of penicillopepsin and neuraminidase crystal structures," <i>J. Mol. Biol.</i> 1994 Oct 14;243(1):100-15.
✓	Kraulis, "MOLSCRIPT: a program to produce both detailed and schematic plots of protein structures," <i>J. Appl. Cryst.</i> 1991 Oct;24:946-950.
✓	Kuntz et al., "A geometric approach to macromolecule-ligand interactions," <i>J. Mol. Biol.</i> 1982 Oct 25;161(2):269-88.
✓	Laemmli, "Cleavage of Structural Proteins during the Assembly of the Head of Bacteriophage T4," <i>Nature</i> , 1970 Aug 15;227(259):680-85.
✓	Laskowski, et al., "PROCHECK: a program to check the stereochemical quality of protein structures," <i>J. Appl Cryst.</i> 1993 Apr;26:283-291
✓	Lattman, "Use of the rotation and translation functions," <i>Methods Enzymol.</i> 1985;115:55-77.
✓	Lauri et al., "CAVEAT: a program to facilitate the design of organic molecules," <i>J Comput Aided Mol Des.</i> 1994 Feb;8(1):51-66.
✓	Martin, "3D database searching in drug design," <i>J. Med. Chem.</i> 1992 Jun 12;35(12):2145-54.
✓	Meng et al., "Automated docking with grid-based energy evaluation," <i>J. Comp. Chem.</i> 1992 May;13(4):505-524.
✓	Merritt et al., "Raster3D Version 2.0. A Program for Photorealistic Molecular Graphics", <i>Acta Crystallogr D Biol Crystallogr.</i> , 1994;50:869-73.
✓	Miranker et al., "Functionality maps of binding sites: a multiple copy simultaneous search method," <i>Proteins: Struct. Funct. Genet.</i> 1991;11(1):29-34.
✓	Moat et al., "Biosynthesis and salvage pathways of pyridine nucleotides. Coenzymes and cofactors, Pyridine Nucleotide Coenzymes," Eds. D. Dolphin et al. John Wiley & Sons, Inc., New York, 1987; vol.II, part B:1-24.
✓	National Institutes of Health, "BLAST 2 Sequences," [online] United States; retrieved October 15, 2001 from the Internet: <URL:http://www.ncbi.nlm.nih.gov/gorf/bl2.html>, 1 pg.
✓	Navaza, "AMoRe: an automated package for molecular replacement," <i>Acta Crystallogr A.</i> 1994 Mar;50:157-163.

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GSAA JAN 08 2002 PATENT & TRADEMARK OFFICE	Nessi et al., "The outB gene of <i>Bacillus subtilis</i> codes for NAD synthetase," <i>J Biol Chem.</i> 1995 Mar 17;270(11):6181-5.
	Nishibata et al., "Automatic creation of drug candidate structures based on receptor structure. Starting point for artificial lead generation," <i>Tetrahedron</i> , 1991;47(43):8985-90.
	Otwinowski, "Maximum likelihood refinement of heavy atom parameters," <i>Isomorphous replacement and anomalous scattering - Proceedings of the CCP4 Study Weekend 25-26 January 1991</i> , (W. Wolf et al., Eds.) Science and Engineering Research Counsel, Daresbury Laboratory, Warrington, U.K. (1991) pp. 80-86.
	Øzment et al., "Structural study of <i>Escherichia coli</i> NAD synthetase: overexpression, purification, crystallization, and preliminary crystallographic analysis," <i>J. Struct. Biol.</i> 1999 Oct;127(3):279-82.
	Ramakrishnan et al., "Crystal structure of globular domain of histone H5 and its implications for nucleosome binding," <i>Nature</i> . 1993 Mar 18;362(6417):219-23.
	Research Collaboratory for Structural Bioinformatics, "Protein Data Bank," [online] United States; retrieved October 15, 2001 from the Internet: <URL: http://www.rcsb.org/pdb/ > 1 page.
	Rizzi et al., "Crystal structure of NH ₃ -dependent NAD ⁺ synthetase from <i>Bacillus subtilis</i> ," <i>EMBO J.</i> 1996 Oct 1;15(19):5125-34.
	Rizzi et al., "Crystallization of NAD ⁺ synthetase from <i>Bacillus subtilis</i> ," <i>Proteins</i> . 1996 Oct;26(2):236-8.
	Rizzi et al., "A novel deamido-NAD ⁺ -binding site revealed by the trapped NAD-adenylate intermediate in the NAD ⁺ synthetase structure," <i>Structure</i> . 1998 Sep 15;6(9):1129-40.
	Rossmann et al., "Chemical and biological evolution of nucleotide-binding protein," <i>Nature</i> . 1974 Jul 19;250(463):194-9.
	Rossmann, ed., <i>The Molecular Replacement Method - A Collection of Papers on the Use of Non-Crystallographic Symmetry</i> , Intl. Sci. Rev. Ser. No. 13, Gordon & Breach, New York, NY; title page, publication page, and table of contents only, 6 pages (1972).
	Sack, "CHAIN - A Crystallographic Modeling Program," <i>J Molecular Graphics</i> . 1988 Dec; 6(4):224-5.

EXAMINER C. Mueller	Date Considered September 14, 2002
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	Sheldrick et al., "Structure solution by iterative peaklist optimization and tangent expansion in space group P1," <i>Acta Crystallogr B</i> . 1995 Aug 1;51(Pt 4):423-31.
	Tatusova et al., "BLAST 2 Sequences, a new tool for comparing protein and nucleotide sequences," <i>FEMS Microbiol Lett</i> . 1999 May 15;174(2):247-50.
	Travis, "Proteins and organic solvents make an eye-opening mix," <i>Science</i> . 1993 Nov 26;262(5138):1374.
	Van Duyne et al., "Atomic structures of the human immunophilin FKBP-12 complexes with FK506 and rapamycin," <i>J Mol Biol</i> . 1993 Jan 5;229(1):105-24.
	White et al., " <i>Biosynthesis of salvage pathways of pyridine nucleotide coenzymes</i> ." Academic Press, New York; title page, publication page, and table of contents only, 13 pages (1982)
	Willison, "An essential gene (efg) located at 38.1 minutes on the Escherichia coli chromosome," <i>J Bacteriol</i> . 1992 Sep;174(17):5765-6.
	Willison et al., "The Escherichia coli efg gene and the Rhodobacter capsulatus adgA gene code for NH3-dependent NAD synthetase," <i>J Bacteriol</i> . 1994 Jun;176(11):3400-2.
	Wyckoff et al., eds., <i>Methods in Enzymology Vol. 114 - Diffraction Methods for Biological Macromolecules</i> , Academic Press, Orlando, FL; title page, publication page, and table of contents only, 5 pages (1985).
	Wyckoff et al., eds., <i>Methods in Enzymology Vol. 115. Diffraction Methods for Biological Macromolecules</i> , Academic Press, Orlando, FL; title page, publication page, and table of contents only, 4 pages (1985).
✓	✓ Zalkin, "NAD synthetase," <i>Methods Enzymol</i> . 1985;113:297-302.

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